Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. - 8. (canceled)

9. (Currently Amended) A pharmaceutical composition comprising:

(i) a compound of formula I

$$O$$
 R^{1}
 $CCH_{2})_{n}$
 R^{2}

or a pharmaceutically acceptable salt, ester, or solvate of the compound, wherein:

n is 1;

X is either O or S;

 R^1 is C_1C_2 - C_9 straight or branched chain alkyl, C_2 - C_9 straight or branched chain alkenyl, aryl, heteroaryl, carbocycle, or heterocycle;

D is a bond, C1-C10 straight or branched chain alkyl, C2-C10 alkenyl or C2-C10 alkynyl;

R² is COOH; carboxylic acid or a carboxylic acid isostere

said alkyl, alkonyl, alkynyl, aryl, heteroaryl, earbocycle, or heterocycle is optionally substituted with one or more substituents selected from hydroxy, halo, haloalkyl, thiocarbonyl, alkoxy, alkenoxy, alkylaryloxy, aryloxy, arylalkyloxy, eyano, nitro, imino, alkylamino, aminoalkyl, sulfhydryl, thioalkyl, alkylthio, sulfonyl, C_1 , C_6 straight or branched chain alkyl, C_2 , C_6 straight or branched chain alkenyl or alkynyl, aryl, aralkyl, heteroaryl, earbocycle, heterocycle, and CO_2R^7 wherein R^7 is hydrogen, C_4 , C_9 straight or branched chain alkyl or C_2 , C_9 straight or branched chain alkenyl; and

(ii) a pharmaceutically acceptable carrier.

Currently Amended) The pharmaceutical composition of claim 9, wherein the compound is selected from the group consisting of compounds of Formula (I) 1, 3, 5, 8, 11, 14, 17, 21, 24-32, 34, 38-40, 44, 45, 47-52, 62, 64-68, 73-98, 101, 102, 106, 108-117 and 119-137 of Tables I, II and III

$$R^{1}$$
 $(CH_{2})_{n}$
 R^{2}
 R^{1}
,wherein,

(a) D is a bond, R₂ is COOH, and

X, n, and R_1 are defined as follows:

No. X n R1

1	Q	1	3,4,5-trimethylphenyl	
3	ō	1	tert-butyl	
<u>5</u>	O	1	cyclopentyl	
8	Q	1	cyclohexyl	
11	O	1	<u>cycloheptyl</u>	
14	O	1	2-thienyl	
17	<u>o</u>	1	2-furyl	
21	<u>o</u>	1	1,1-dimethylpentyl	

<u>or</u>

(b) n, X, D, R₂, and R₁ are defined as follows:

No.	n	X	D	<u>R₂</u>	<u>R</u> 1
83	1	O	bond	СООН	<u>α-Methylphenyl</u>
84	1	Ō	bond	СООН	4-Methylphenyl
<u>137</u>	1	<u>o</u>	bond	СООН	1,1-dimethylpropyl

14. – 82. (Cancelled)

$$R^{1}$$
 $(CH_{2})_{n}$
 R^{2}

or a pharmaceutically acceptable salt, ester, or solvate of the compound, wherein:

X is O;

n is 1;

D is a bond;

 R^2 is – COOH;

R¹ is 1,1-dimethylpropyl, and

(ii) a pharmaceutically acceptable carrier.